Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111115\
Data File : BE091149.D
Acq On : 12 Nov 2015 18:18
Operator : UM/NP
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTD0.496

Quant Time: Nov 13 17:26:19 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE102715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 09 05:49:45 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	87	-0.06
2	SURR	1,4-Dioxane-d8	0.288	0.242	16.0	69	-0.22
3		1,4-Dioxane	0.325	0.271	16.6	71	-0.21
4	I	Naphthalene-d8	1.000	1.000	0.0	99	-0.04
5		Naphthalene	0.986	0.950	3.7	92	-0.04
6	SURR	2-Methylnaphthalene-d10	0.546	0.552	-1.1	99	-0.05
7		2-Methylnaphthalene	0.643	0.648	-0.8	98	-0.05
8	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
9		Acenaphthylene	1.998	1.956	2.1	100	0.00
10	C	Acenaphthene	1.368	1.303	4.8	98	0.00
11		Fluorene	1.757	1.586	9.7	89	0.00
12	I	Phenanthrene-d10	1.000	1.000	0.0	84	0.02
13		Pentachlorophenol	0.038	0.055	-44.7#	150#	0.01
14		Phenanthrene	4.619	4.405	4.6	79	0.02
15		Anthracene	4.200	4.299	-2.4	89	0.01
16	SURR	Fluoranthene-d10	1.127	1.030	8.6	82	0.03
17	C	Fluoranthene	5.406	4.924	8.9	80	0.03
18	I	Chrysene-d12	1.000	1.000	0.0	82	0.04
19		Pyrene	4.544	4.524	0.4	81	0.03
20		Benzo(a)anthracene	1.027	1.058	-3.0	87	0.04
21		Chrysene	1.183	1.111	6.1	76	0.04
22	I	Perylene-d12	1.000	1.000	0.0	93	0.05
23		Benzo(b)fluoranthene	1.249	1.164	6.8	85	0.05
24		Benzo(k)fluoranthene	1.281	1.126	12.1	78	0.05
25	C	Benzo(a)pyrene	1.150	1.108	3.7	89	0.05
26		Indeno(1,2,3-cd)pyrene	4.847	4.380	9.6	82	0.09
27		Dibenzo(a,h)anthracene	1.155	1.114	3.5	90	0.09
28		Benzo(g,h,i)perylene	5.083	4.601	9.5	84	0.09

(#) = Out of Range SPCC's out = 0 CCC's out = 0